



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:18 am GMT

PDB ID : 1AB8  
Title : RAT TYPE II ADENYLYL CYCLASE C2 DOMAIN/FORSKOLIN COM-  
PLEX  
Authors : Zhang, G.; Hurley, J.H.  
Deposited on : 1997-02-04  
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

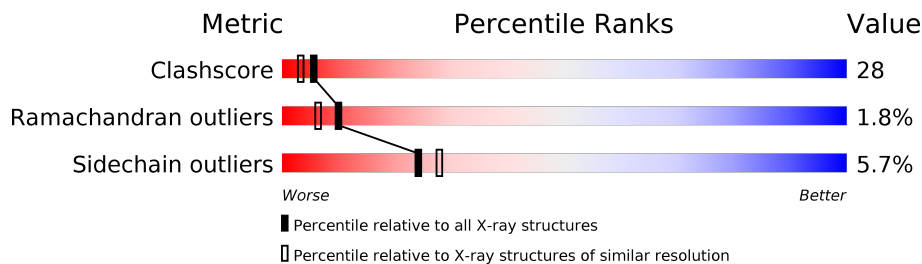
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	220	
1	B	220	

## 2 Entry composition [i](#)

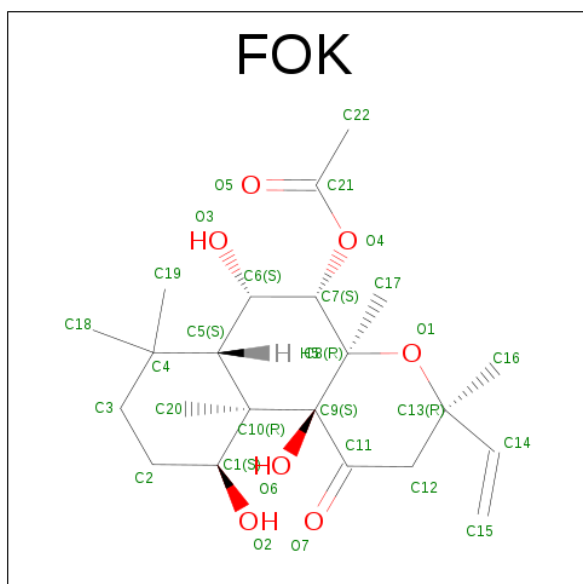
There are 3 unique types of molecules in this entry. The entry contains 2885 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ADENYLYL CYCLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	0	0
			1355	870	215	261	9			
1	B	176	Total	C	N	O	S	0	0	0
			1359	869	218	263	9			

- Molecule 2 is FORSKOLIN (three-letter code: FOK) (formula:  $C_{22}H_{34}O_7$ ).



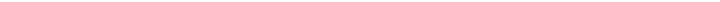
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			29	22	7		
2	B	1	Total	C	O	0	0
			29	22	7		

- Molecule 3 is water.

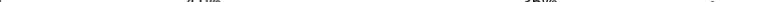
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	62	Total 62	O 62	0	0
3	B	51	Total 51	O 51	0	0

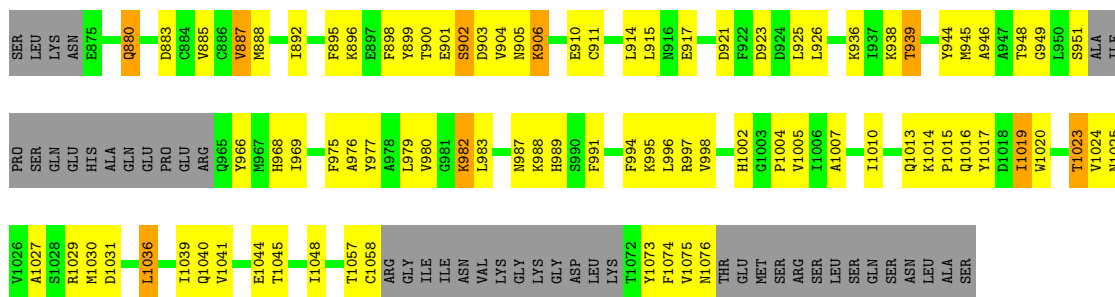


Note EDS was not executed.

Chain A:  49% 28% 3% 20%



Chain B:  41% 35% 5% 19%



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.30Å 81.30Å 180.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20	Depositor
% Data completeness (in resolution range)	54.2 (6.00-2.20)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, $R_{free}$	0.219 , 0.284	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2885	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FOK

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.56	1/1380 (0.1%)	1.03	2/1869 (0.1%)
1	B	0.49	0/1383	0.72	1/1871 (0.1%)
All	All	0.53	1/2763 (0.0%)	0.89	3/3740 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1056	CYS	C-N	7.75	1.51	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1056	CYS	O-C-N	-31.29	72.63	122.70
1	B	1019	ILE	N-CA-C	-5.68	95.66	111.00
1	A	1019	ILE	N-CA-C	-5.33	96.62	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1056	CYS	Mainchain

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1355	0	1326	63	0
1	B	1359	0	1328	84	0
2	A	29	0	33	11	0
2	B	29	0	33	14	0
3	A	62	0	0	10	0
3	B	51	0	0	19	0
All	All	2885	0	2720	156	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 28.

All (156) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1058:CYS:HG	1:B:1073:TYR:HE1	1.03	1.00
1:A:1018:ASP:HB2	3:A:80:HOH:O	1.61	0.99
2:A:1:FOK:H201	2:A:1:FOK:H173	1.46	0.97
1:B:1024:VAL:HG23	3:B:48:HOH:O	1.65	0.97
2:B:2:FOK:H173	2:B:2:FOK:H201	1.47	0.95
1:A:997:ARG:HD3	1:A:1031:ASP:O	1.76	0.85
1:A:1034:GLY:HA3	3:A:81:HOH:O	1.77	0.84
1:B:899:TYR:N	3:B:49:HOH:O	2.11	0.82
1:B:997:ARG:HD3	1:B:1031:ASP:O	1.80	0.81
1:A:888:MET:HE1	1:A:979:LEU:HD11	1.61	0.81
1:B:1024:VAL:N	3:B:48:HOH:O	2.15	0.80
2:A:1:FOK:H202	2:A:1:FOK:H193	1.63	0.79
1:B:1005:VAL:HG11	1:B:1019:ILE:HG23	1.67	0.76
1:B:946:ALA:O	3:B:44:HOH:O	2.04	0.76
1:A:1005:VAL:HG11	1:A:1019:ILE:HG23	1.66	0.76
1:B:945:MET:HE1	2:B:2:FOK:H32	1.71	0.72
1:B:1023:THR:HG23	3:B:47:HOH:O	1.88	0.72
1:A:1041:VAL:HG22	1:A:1045:THR:HB	1.72	0.72
1:A:1019:ILE:HG22	1:A:1024:VAL:HG21	1.72	0.72
1:B:1041:VAL:HG22	1:B:1045:THR:HB	1.72	0.71
1:A:902:SER:H	1:A:905:ASN:HB3	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2:FOK:H193	2:B:2:FOK:H202	1.75	0.68
1:A:1023:THR:HG22	3:A:84:HOH:O	1.93	0.67
1:A:995:LYS:HB3	1:A:1036:LEU:HD12	1.77	0.66
1:B:991:PHE:HB2	3:B:35:HOH:O	1.95	0.66
1:A:877:LEU:HA	3:A:89:HOH:O	1.95	0.65
1:B:905:ASN:O	1:B:906:LYS:HB2	1.97	0.65
1:B:1005:VAL:CG1	1:B:1019:ILE:HG23	2.27	0.65
1:B:905:ASN:O	1:B:910:GLU:HG2	1.97	0.65
1:B:995:LYS:HB3	1:B:1036:LEU:HD12	1.78	0.65
1:B:899:TYR:HB3	3:B:49:HOH:O	1.97	0.65
1:A:969:ILE:HD12	1:A:1048:ILE:HG21	1.79	0.65
2:A:1:FOK:H172	1:B:896:LYS:NZ	2.11	0.65
1:B:892:ILE:N	3:B:40:HOH:O	2.17	0.65
1:B:1058:CYS:SG	1:B:1073:TYR:CE1	2.90	0.64
1:B:969:ILE:HD12	1:B:1048:ILE:HG21	1.81	0.63
1:B:914:LEU:HD21	1:B:994:PHE:HE2	1.63	0.63
1:A:1018:ASP:CB	3:A:80:HOH:O	2.34	0.62
1:B:1058:CYS:SG	1:B:1073:TYR:HE1	2.17	0.62
1:A:1005:VAL:CG1	1:A:1019:ILE:HG23	2.29	0.62
1:A:945:MET:HE1	2:A:1:FOK:H32	1.81	0.62
2:B:2:FOK:C17	2:B:2:FOK:H201	2.26	0.62
1:B:880:GLN:HA	3:B:76:HOH:O	2.00	0.62
1:A:995:LYS:CB	1:A:1036:LEU:HD12	2.31	0.61
2:A:1:FOK:H201	2:A:1:FOK:C17	2.25	0.61
1:B:1019:ILE:HG22	1:B:1024:VAL:HG21	1.80	0.61
1:B:995:LYS:CB	1:B:1036:LEU:HD12	2.30	0.61
1:A:896:LYS:NZ	2:B:2:FOK:H172	2.16	0.60
1:B:1024:VAL:HG11	2:B:2:FOK:H1	1.84	0.59
1:B:888:MET:O	3:B:44:HOH:O	2.16	0.59
1:A:1019:ILE:HG22	1:A:1024:VAL:CG2	2.32	0.59
1:A:1057:THR:HG22	1:A:1058:CYS:N	2.18	0.58
1:A:1018:ASP:O	1:A:1019:ILE:HD13	2.04	0.58
1:A:988:LYS:HE2	3:A:119:HOH:O	2.04	0.58
1:B:948:THR:HB	1:B:968:HIS:HB2	1.86	0.58
1:A:914:LEU:HD21	1:A:994:PHE:HE2	1.69	0.57
1:A:889:PHE:HE1	1:A:945:MET:HE2	1.70	0.57
1:B:1057:THR:O	1:B:1073:TYR:HA	2.05	0.56
1:A:1049:LEU:HD13	1:A:1075:VAL:HG21	1.87	0.56
1:A:906:LYS:HB2	1:A:910:GLU:HG2	1.88	0.56
1:B:887:VAL:HG22	1:B:1027:ALA:CB	2.36	0.55
1:A:1057:THR:HG22	1:A:1058:CYS:H	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:883:ASP:HA	1:B:1004:PRO:HA	1.89	0.55
1:A:896:LYS:HZ2	2:B:2:FOK:H172	1.71	0.54
1:B:899:TYR:CE1	1:B:901:GLU:HG3	2.42	0.54
1:B:1025:ASN:HB2	2:B:2:FOK:H171	1.88	0.54
1:B:1019:ILE:HG22	1:B:1024:VAL:CG2	2.37	0.53
1:A:906:LYS:HD2	1:A:910:GLU:OE2	2.08	0.53
2:A:1:FOK:H121	2:A:1:FOK:H171	1.89	0.53
1:B:901:GLU:O	1:B:902:SER:HB2	2.09	0.53
1:A:889:PHE:CE1	1:A:945:MET:HE2	2.42	0.53
1:A:953:ILE:O	1:A:953:ILE:HG23	2.09	0.52
1:B:1024:VAL:HG12	2:B:2:FOK:H203	1.92	0.52
1:A:976:ALA:O	1:A:980:VAL:HG23	2.10	0.51
1:A:1075:VAL:HG12	1:A:1076:ASN:N	2.26	0.51
1:B:976:ALA:O	1:B:980:VAL:HG23	2.11	0.51
1:B:925:LEU:HD12	1:B:979:LEU:HD23	1.93	0.50
2:B:2:FOK:H171	2:B:2:FOK:H121	1.93	0.50
1:A:1054:TYR:HB3	1:A:1075:VAL:HG13	1.93	0.50
1:B:900:THR:OG1	1:B:905:ASN:ND2	2.45	0.50
2:A:1:FOK:H172	1:B:896:LYS:HZ3	1.76	0.50
1:A:948:THR:HB	1:A:968:HIS:HB2	1.93	0.50
1:B:1014:LYS:N	1:B:1015:PRO:HD3	2.26	0.50
1:A:1024:VAL:HG11	2:A:1:FOK:H1	1.95	0.49
1:B:1007:ALA:HB1	1:B:1017:TYR:CZ	2.48	0.49
1:B:895:PHE:CZ	1:B:915:LEU:HB2	2.48	0.49
1:A:887:VAL:HG22	1:A:1027:ALA:CB	2.44	0.48
1:A:942:SER:C	3:A:79:HOH:O	2.52	0.48
1:A:978:ALA:O	3:A:126:HOH:O	2.20	0.48
1:B:880:GLN:HG3	3:B:87:HOH:O	2.14	0.48
1:B:939:THR:HG22	1:B:944:TYR:HA	1.96	0.47
1:B:946:ALA:N	3:B:44:HOH:O	2.17	0.47
1:B:1030:MET:CE	1:B:1040:GLN:HG2	2.43	0.47
1:B:888:MET:HE1	1:B:975:PHE:CE2	2.50	0.47
1:A:1016:GLN:NE2	1:B:938:LYS:HG2	2.30	0.47
1:B:1005:VAL:HG22	3:B:48:HOH:O	2.14	0.46
1:A:1004:PRO:O	1:A:1023:THR:HG23	2.15	0.46
1:B:899:TYR:CA	3:B:49:HOH:O	2.59	0.46
2:A:1:FOK:H172	1:B:896:LYS:HZ2	1.81	0.46
1:B:921:ASP:HB3	1:B:982:LYS:HG2	1.98	0.46
1:B:906:LYS:HD2	1:B:910:GLU:OE2	2.16	0.45
1:B:977:TYR:OH	1:B:1075:VAL:HG22	2.17	0.45
1:A:966:TYR:HB3	1:A:1048:ILE:CG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:987:ASN:HD21	1:B:994:PHE:HB2	1.81	0.45
1:A:1018:ASP:CG	3:A:80:HOH:O	2.55	0.44
1:B:899:TYR:HA	1:B:911:CYS:SG	2.58	0.44
1:A:1016:GLN:NE2	1:B:939:THR:H	2.15	0.44
1:A:1049:LEU:HB3	1:A:1054:TYR:HB2	1.99	0.44
1:B:1024:VAL:HG12	2:B:2:FOK:C20	2.48	0.44
1:B:996:LEU:HA	1:B:996:LEU:HD23	1.75	0.44
1:A:886:CYS:HB3	1:A:948:THR:OG1	2.18	0.44
1:B:1004:PRO:O	1:B:1023:THR:HG23	2.17	0.44
1:B:899:TYR:CB	3:B:49:HOH:O	2.59	0.44
1:B:898:PHE:O	1:B:900:THR:HG23	2.18	0.44
1:A:883:ASP:HA	1:A:1004:PRO:HA	2.00	0.43
2:B:2:FOK:H173	2:B:2:FOK:C20	2.34	0.43
1:B:925:LEU:CD1	1:B:979:LEU:HD23	2.48	0.43
1:A:1041:VAL:CG2	1:A:1045:THR:HB	2.47	0.43
1:B:1025:ASN:HD21	1:B:1029:ARG:HH21	1.65	0.43
1:A:974:GLU:HG3	3:A:135:HOH:O	2.17	0.43
2:A:1:FOK:C12	2:A:1:FOK:H171	2.49	0.43
1:A:900:THR:O	1:A:905:ASN:OD1	2.37	0.43
1:B:1010:ILE:HD13	1:B:1020:TRP:CZ2	2.54	0.43
1:A:1013:GLN:HG2	1:B:923:ASP:HB3	2.01	0.43
1:A:1006:ILE:HG13	3:B:90:HOH:O	2.18	0.43
1:A:1002:HIS:HE1	1:A:1044:GLU:HG2	1.84	0.42
1:B:1002:HIS:HE1	1:B:1044:GLU:HG2	1.85	0.42
1:B:888:MET:HG3	1:B:998:VAL:HG13	2.01	0.42
1:A:966:TYR:HB3	1:A:1048:ILE:HG23	2.00	0.42
1:B:892:ILE:HB	3:B:40:HOH:O	2.18	0.42
1:B:926:LEU:HD11	1:B:936:LYS:HB2	2.02	0.42
1:B:914:LEU:O	1:B:917:GLU:HB3	2.19	0.42
1:B:1030:MET:HE1	1:B:1040:GLN:HG2	2.02	0.42
1:B:949:GLY:C	1:B:951:SER:H	2.21	0.42
2:B:2:FOK:C17	2:B:2:FOK:C20	2.97	0.42
1:B:888:MET:HE1	1:B:975:PHE:HE2	1.82	0.42
1:A:988:LYS:HG3	1:A:989:HIS:CD2	2.55	0.42
1:A:898:PHE:HE2	1:A:992:ASN:HD21	1.68	0.42
1:B:1039:ILE:O	1:B:1074:PHE:HA	2.20	0.42
1:B:1041:VAL:CG2	1:B:1045:THR:HB	2.46	0.42
1:A:938:LYS:HE3	1:A:945:MET:HE1	2.02	0.41
1:A:1039:ILE:O	1:A:1074:PHE:HA	2.20	0.41
1:B:892:ILE:CA	3:B:40:HOH:O	2.66	0.41
1:A:987:ASN:HD21	1:A:994:PHE:HB2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:888:MET:HB3	1:A:972:MET:CE	2.50	0.41
1:A:1016:GLN:HE21	1:B:938:LYS:HG2	1.85	0.41
1:A:925:LEU:HD12	1:A:979:LEU:HD23	2.03	0.41
1:B:946:ALA:C	3:B:44:HOH:O	2.56	0.41
1:B:914:LEU:HD21	1:B:994:PHE:CE2	2.50	0.41
1:B:988:LYS:HG3	1:B:989:HIS:CD2	2.55	0.41
1:A:1010:ILE:HD13	1:A:1020:TRP:CZ2	2.56	0.41
1:A:983:LEU:HA	1:A:983:LEU:HD23	1.80	0.41
1:B:885:VAL:O	1:B:1002:HIS:HA	2.21	0.41
2:B:2:FOK:C12	2:B:2:FOK:H171	2.52	0.40
2:A:1:FOK:C20	2:A:1:FOK:C17	2.96	0.40
1:B:902:SER:O	1:B:904:VAL:N	2.55	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	171/220 (78%)	158 (92%)	11 (6%)	2 (1%)	15	12
1	B	170/220 (77%)	153 (90%)	13 (8%)	4 (2%)	7	4
All	All	341/440 (78%)	311 (91%)	24 (7%)	6 (2%)	10	6

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	954	PRO
1	B	903	ASP
1	A	1057	THR
1	B	902	SER
1	B	966	TYR
1	B	906	LYS

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	148/191 (78%)	141 (95%)	7 (5%)	30	37
1	B	149/191 (78%)	139 (93%)	10 (7%)	19	21
All	All	297/382 (78%)	280 (94%)	17 (6%)	24	28

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	887	VAL
1	A	954	PRO
1	A	982	LYS
1	A	983	LEU
1	A	1014	LYS
1	A	1023	THR
1	A	1036	LEU
1	B	880	GLN
1	B	887	VAL
1	B	939	THR
1	B	982	LYS
1	B	983	LEU
1	B	1013	GLN
1	B	1016	GLN
1	B	1023	THR
1	B	1036	LEU
1	B	1076	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	987	ASN
1	A	989	HIS
1	A	992	ASN
1	A	1001	ASN
1	A	1040	GLN

*Continued on next page...*

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Mol	Chain	Res	Type
1	B	880	GLN
1	B	905	ASN
1	B	968	HIS
1	B	987	ASN
1	B	989	HIS
1	B	992	ASN
1	B	1001	ASN
1	B	1013	GLN
1	B	1016	GLN
1	B	1040	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	FOK	A	1	-	26,31,31	1.43	5 (19%)	31,54,54	1.38	4 (12%)
2	FOK	B	2	-	26,31,31	1.48	7 (26%)	31,54,54	1.48	7 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FOK	A	1	-	-	0/7/80/80	0/3/3/3
2	FOK	B	2	-	-	0/7/80/80	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	FOK	C15-C14	-2.42	1.15	1.29
2	B	2	FOK	C15-C14	-2.37	1.15	1.29
2	A	1	FOK	O6-C9	2.05	1.48	1.42
2	A	1	FOK	C22-C21	2.10	1.56	1.49
2	B	2	FOK	O6-C9	2.12	1.48	1.42
2	B	2	FOK	C18-C4	2.13	1.58	1.53
2	B	2	FOK	C22-C21	2.18	1.57	1.49
2	B	2	FOK	C19-C4	2.20	1.58	1.53
2	A	1	FOK	C17-C8	2.80	1.56	1.51
2	B	2	FOK	C17-C8	2.81	1.56	1.51
2	B	2	FOK	C3-C4	3.80	1.62	1.54
2	A	1	FOK	C3-C4	3.93	1.62	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	FOK	C2-C3-C4	-3.19	107.89	113.42
2	B	2	FOK	C2-C3-C4	-3.05	108.13	113.42
2	B	2	FOK	C7-O4-C21	-2.49	114.18	117.81
2	B	2	FOK	C2-C1-C10	-2.34	108.32	112.06
2	A	1	FOK	C2-C1-C10	-2.24	108.49	112.06
2	B	2	FOK	C19-C4-C3	-2.16	104.46	109.00
2	B	2	FOK	C10-C5-C4	2.02	118.16	116.48
2	A	1	FOK	C13-C14-C15	2.25	138.74	128.76
2	B	2	FOK	C13-C14-C15	2.35	139.20	128.76
2	B	2	FOK	C3-C2-C1	2.72	115.67	111.47
2	A	1	FOK	C3-C2-C1	3.30	116.56	111.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	FOK	11	0
2	B	2	FOK	14	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.